

Data Path : U:\HPCHEM1\BNA F\DATA\BF101915\
 Data File : BF082352.D
 Acq On : 19 Oct 2015 13:18
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

MMdadoda
 10/20/2015 6:25:41 PM

Quant Time: Oct 20 01:57:44 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 01:46:21 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	149596	20.00	ng	-0.03
21) Naphthalene-d8	8.09	136	587997	20.00	ng	-0.03
38) Acenaphthene-d10	9.85	164	295216	20.00	ng	-0.03
63) Phenanthrene-d10	11.35	188	587022	20.00	ng	-0.02
75) Chrysene-d12	14.09	240	509909	20.00	ng	0.08
86) Perylene-d12	15.67	264	402116	20.00	ng	0.23

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	625160	84.34	ng	-0.03
7) Phenol-d6	6.45	99	783884	81.27	ng	-0.03
23) Nitrobenzene-d5	7.38	82	681127	77.79	ng	-0.03
41) 2,4,6-Tribromophenol	10.65	330	185828	67.12	ng	-0.02
44) 2-Fluorobiphenyl	9.18	172	1259721	70.76	ng	-0.03
78) Terphenyl-d14	12.95	244	1303516	67.74	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.34	88	140970	37.85	ng	97
3) Pyridine	3.05	79	383468	44.27	ng	97
4) n-Nitrosodimethylamine	3.03	42	149857	40.79	ng	97
6) Aniline	6.47	93	497991	40.04	ng	# 49
8) 2-Chlorophenol	6.59	128	352701	38.98	ng	98
9) Benzaldehyde	6.35	77	210636	42.76	ng	98
10) Phenol	6.47	94	440044	39.57	ng	# 71
11) bis(2-Chloroethyl)ether	6.55	93	353732	37.61	ng	96
12) 1,3-Dichlorobenzene	6.74	146	418581	39.72	ng	99
13) 1,4-Dichlorobenzene	6.82	146	431620	39.22	ng	99
14) 1,2-Dichlorobenzene	6.97	146	355526	34.02	ng	99
15) Benzyl Alcohol	6.96	79	279083	41.91	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.09	45	461825	35.26	ng	89
17) 2-Methylphenol	7.07	107	276164	38.59	ng	98
18) Hexachloroethane	7.31	117	145235	38.56	ng	95
19) n-Nitroso-di-n-propylamine	7.23	70	236550	34.92	ng	# 91
20) 3+4-Methylphenols	7.23	107	346057	40.59	ng	# 78
22) Acetophenone	7.22	105	459342	39.49	ng	# 95
24) Nitrobenzene	7.41	77	385200	40.64	ng	# 83
25) Isophorone	7.65	82	702914	39.78	ng	95
26) 2-Nitrophenol	7.71	139	194012	41.00	ng	# 88
27) 2,4-Dimethylphenol	7.76	122	306751	40.23	ng	98
28) bis(2-Chloroethoxy)methane	7.85	93	388708	37.80	ng	98
29) 2,4-Dichlorophenol	7.97	162	325223	42.67	ng	96
30) 1,2,4-Trichlorobenzene	8.03	180	335847	38.23	ng	97
31) Naphthalene	8.11	128	978056	38.37	ng	99
32) Benzoic acid	7.91	122	158350	30.95	ng	98
33) 4-Chloroaniline	8.17	127	412444	38.96	ng	95
34) Hexachlorobutadiene	8.23	225	195855	35.78	ng	97
35) Caprolactam	8.57	113	86457	39.09	ng	# 83
36) 4-Chloro-3-methylphenol	8.66	107	295158	39.60	ng	99
37) 2-Methylnaphthalene	8.81	142	678948	38.42	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	344970	39.29	ng	97
40) Hexachlorocyclopentadiene	8.96	237	117761	32.06	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.10	196	242990	41.66	nq	98
43) 2,4,5-Trichlorophenol	9.14	196	242896	38.45	nq	98
45) 1,1'-Biphenyl	9.28	154	857188	38.08	nq	97
46) 2-Chloronaphthalene	9.30	162	638239	36.01	nq	# 94
47) 2-Nitroaniline	9.41	65	199678	41.04	nq	93
48) Acenaphthylene	9.71	152	1023133	37.06	nq	100
49) Dimethylphthalate	9.59	163	779170	37.48	nq	100
50) 2,6-Dinitrotoluene	9.66	165	195804	44.64	nq	91
51) Acenaphthene	9.90	154	613264	37.00	nq	99
52) 3-Nitroaniline	9.83	138	209308	42.24	nq	94
53) 2,4-Dinitrophenol	9.93	184	87243	39.14	nq	98
54) Dibenzofuran	10.07	168	880280	38.69	nq	93
55) 4-Nitrophenol	9.99	139	115385	41.58	nq	# 83
56) 2,4-Dinitrotoluene	10.06	165	224390	40.93	nq	# 90
57) Fluorene	10.41	166	744486	39.84	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.18	232	195642	37.90	nq	97
59) Diethylphthalate	10.29	149	812044	41.91	nq	99
60) 4-Chlorophenyl-phenylether	10.40	204	360948	42.16	nq	90
61) 4-Nitroaniline	10.45	138	200872	41.80	nq	96
62) Azobenzene	10.56	77	776544	40.95	nq	93
64) 4,6-Dinitro-2-methylphenol	10.47	198	130812	39.71	nq	83
65) n-Nitrosodiphenylamine	10.53	169	706205	40.18	nq	99
66) 4-Bromophenyl-phenylether	10.89	248	221589	37.72	nq	# 92
67) Hexachlorobenzene	10.95	284	213646	33.63	nq	# 65
68) Atrazine	11.05	200	190531	34.22	nq	98
69) Pentachlorophenol	11.15	266	128359	39.26	nq	99
70) Phenanthrene	11.37	178	1123376	39.04	nq	99
71) Anthracene	11.42	178	1100252	37.11	nq	99
72) Carbazole	11.58	167	984003	36.38	nq	100
73) Di-n-butylphthalate	11.91	149	1220772	36.59	nq	99
74) Fluoranthene	12.56	202	1105613	35.78	nq	97
76) Benzidine	12.70	184	503019	34.04	nq	99
77) Pyrene	12.80	202	1209241	39.96	nq	99
79) Butylbenzylphthalate	13.45	149	548403	39.71	nq	99
80) Benzo(a)anthracene	14.07	228	973290	36.69	nq	99
81) 3,3'-Dichlorobenzidine	14.05	252	408924	40.61	nq	# 96
82) Chrysene	14.12	228	1006027	35.99	nq	99
83) Bis(2-ethylhexyl)phthalate	14.07	149	744654	37.80	nq	100
84) Di-n-octyl phthalate	14.79	149	1330998	39.34	nq	99
85) Indeno(1,2,3-cd)pyrene	17.12	276	829554	37.33	nq	98
87) Benzo(b)fluoranthene	15.25	252	1023345	41.83	nq	100
88) Benzo(k)fluoranthene	15.28	252	683844m	33.26	nq	
89) Benzo(a)pyrene	15.61	252	792007	37.67	nq	100
90) Dibenzo(a,h)anthracene	17.13	278	705167	40.93	nq	99
91) Benzo(g,h,i)perylene	17.54	276	687426	41.07	nq	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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